EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	5229	((568/667) or (568/669) or (568/670) or (526/72) or (528/401) or (528/402) or (430/311)).CCLS.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	OFF	2006/04/15 12:30
L2	7	norborn\$6 and (hexafluoropropanol or trifluoromethyl near2 carbinol)	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/04/15 12:40
L3	0	l1 and l2	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/04/15 12:32
L4	7536	carr.in. or markley.in. or abdourazak. in. or marsella.in.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/04/15 12:41
L5	44228	carbocyclic	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR.	ON	2006/04/15 12:41
L6	27	I4 and I5	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/04/15 12:46

Connecting via Winsock to STN

```
Welcome to STN International! Enter x:x
```

LOGINID:SSSPTA1204rxw

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

```
* * * * * * * * * * Welcome to STN International   * * * * * * * * *
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NEWS
                Web Page URLs for STN Seminar Schedule - N. America
NEWS
                 "Ask CAS" for self-help around the clock
NEWS
        DEC 23
                New IPC8 SEARCH, DISPLAY, and SELECT fields in USPATFULL/
                USPAT2
NEWS
         JAN 13
                IPC 8 searching in IFIPAT, IFIUDB, and IFICDB
NEWS
        JAN 13
                New IPC 8 SEARCH, DISPLAY, and SELECT enhancements added to
                 INPADOC
NEWS 6
                Pre-1988 INPI data added to MARPAT
        JAN 17
NEWS
        JAN 17
                IPC 8 in the WPI family of databases including WPIFV
NEWS 8
        JAN 30
                Saved answer limit increased
NEWS 9
        FEB 21
                STN AnaVist, Version 1.1, lets you share your STN AnaVist
                visualization results
NEWS 10
        FEB 22
                The IPC thesaurus added to additional patent databases on STN
NEWS 11 FEB 22
                Updates in EPFULL; IPC 8 enhancements added
NEWS 12 FEB 27
                New STN AnaVist pricing effective March 1, 2006
NEWS 13 FEB 28 MEDLINE/LMEDLINE reload improves functionality
NEWS 14 FEB 28
                TOXCENTER reloaded with enhancements
NEWS 15 FEB 28 REGISTRY/ZREGISTRY enhanced with more experimental spectral
                property data
NEWS 16 MAR 01
                INSPEC reloaded and enhanced
NEWS 17 MAR 03
                Updates in PATDPA; addition of IPC 8 data without attributes
NEWS 18 MAR 08 X.25 communication option no longer available after June 2006
NEWS 19 MAR 22 EMBASE is now updated on a daily basis
NEWS 20 APR 03 New IPC 8 fields and IPC thesaurus added to PATDPAFULL
NEWS 21 APR 03 Bibliographic data updates resume; new IPC 8 fields and IPC
                thesaurus added in PCTFULL
                STN AnaVist $500 visualization usage credit offered
NEWS 22 APR 04
NEWS 23 APR 12 LINSPEC, learning database for INSPEC, reloaded and enhanced
NEWS 24 APR 12
                Improved structure highlighting in FQHIT and QHIT display
                in MARPAT
NEWS 25 APR 12
                Derwent World Patents Index to be reloaded and enhanced during
                second quarter; strategies may be affected
```

NEWS EXPRESS FEBRUARY 15 CURRENT VERSION FOR WINDOWS IS V8.01a,
CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
AND CURRENT DISCOVER FILE IS DATED 19 DECEMBER 2005.
V8.0 AND V8.01 USERS CAN OBTAIN THE UPGRADE TO V8.01a AT http://download.cas.org/express/v8.0-Discover/

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FILE 'HOME' ENTERED AT 13:08:37 ON 15 APR 2006

=> file reg
COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.21 0.21

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 13:08:53 ON 15 APR 2006 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2006 American Chemical Society (ACS)

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STRUCTURE FILE UPDATES: 13 APR 2006 HIGHEST RN 880388-58-9 DICTIONARY FILE UPDATES: 13 APR 2006 HIGHEST RN 880388-58-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH January 6, 2006

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Structure search iteration limits have been increased. See ${\tt HELP\ SLIMITS}$ for details.

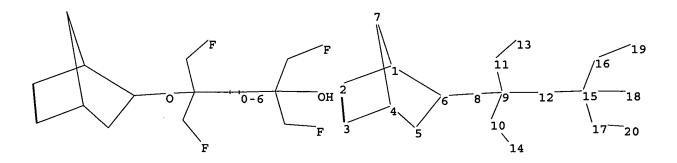
REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/ONLINE/UG/regprops.html

=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=>
Uploading C:\Documents and Settings\rkeys\My
Documents\STNEXP4\TEMPLATE\STANDARD\10784377.str



chain nodes :
8 9 10 11 12 13 14 15 16 17 18 19 20
ring nodes :
1 2 3 4 5 6 7
chain bonds :
6-8 8-9 9-10 9-11 9-12 10-14 11-13 12-15 15-16 15-17 15-18 16-19 17-20
ring bonds :
1-2 1-6 1-7 2-3 3-4 4-5 4-7 5-6
exact/norm bonds :
6-8 8-9 15-18
exact bonds :
1-2 1-6 1-7 2-3 3-4 4-5 4-7 5-6 9-10 9-11 9-12 10-14 11-13 12-15
15-16 15-17 16-19 17-20
isolated ring systems :
containing 1 :

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:CLASS 9:CLASS 10:CLASS 11:CLASS 12:CLASS 13:CLASS 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS

L1 STRUCTURE UPLOADED

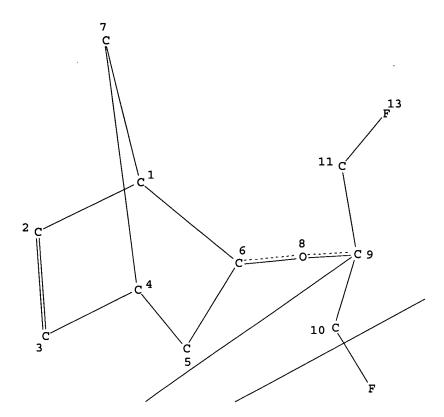
=> que L1

L2 QUE L1

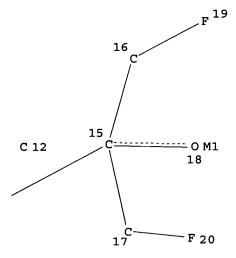
=> d

L2 HAS NO ANSWERS

L1 STR



Page 1-A



Page 1-B

NSPEC IS C ΑT 8 NSPEC IS C 9 AT NSPEC IS C ΑT 10 NSPEC IS C AT11 NSPEC IS C AΤ 12 NSPEC IS C 13 AΤ NSPEC IS C AΤ 14 NSPEC IS C ΑT 15 NSPEC IS C AT 16 NSPEC IS C ΑT 17 NSPEC IS C AT 18 NSPEC IS C TA19 NSPEC IS C AΤ 20 NSPEC IS C AΤ 21 DEFAULT MLEVEL IS ATOM

MLEVEL IS CLASS AT 8 9 10 11 12 13 14 15 16 17 18 19 20 DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RSPEC I

NUMBER OF NODES IS 21

STEREO ATTRIBUTES: NONE L2 QUE L1

=> s 12

SAMPLE SEARCH INITIATED 13:09:19 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 2 TO ITERATE

100.0% PROCESSED SEARCH TIME: 00.00.01 2 ITERATIONS

1 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 2 TO 124
PROJECTED ANSWERS: 1 TO 80

L3 1 SEA SSS SAM L1

=> d scan

L3 1 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN

IN 3-Nonanol, 7-(bicyclo[2.2.1]hept-5-en-2-yloxy)-1,1,1,2,2,8,8,9,9decafluoro-5-methyl-3-[1,2,2,2-tetrafluoro-1-(trifluoromethyl)ethyl]-7(trifluoromethyl)- (9CI)

MF C21 H18 F20 O2

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> 0

0 IS NOT A RECOGNIZED COMMAND

The previous command name entered was not recognized by the system. For a list of commands available to you in the current file, enter "HELP COMMANDS" at an arrow prompt (=>).

=> s 12 ful

FULL SEARCH INITIATED 13:09:58 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 55 TO ITERATE

100.0% PROCESSED 55 ITERATIONS 34 ANSWERS

SEARCH TIME: 00.00.01

L4 34 SEA SSS FUL L1

=> d scan

L4 34 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN

IN 2,9-Decanediol, 4-[2-(bicyclo[2.2.1]hept-5-en-2-yloxy)-3,3,3-trifluoro-2 (trifluoromethyl)propyl]-1,1,1,10,10,10-hexafluoro-7-methyl-2,9bis(trifluoromethyl)-, polymer with 4-[5-(bicyclo[2.2.1]hept-5-en-2-yloxy)-6,6,6-trifluoro-3-methyl-5-(trifluoromethyl)hexyl]-1,1,1,7,7,7-hexafluoro-2,6-bis(trifluoromethyl)-2,6-heptanediol (9CI)

MF (C24 H26 F18 O3 . C24 H26 F18 O3)x

CI PMS

CM 1

CM 2

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> file caplus
COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 167.38 167.59

FULL ESTIMATED COST

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FILE COVERS 1907 - 15 Apr 2006 VOL 144 ISS 17 FILE LAST UPDATED: 14 Apr 2006 (20060414/ED)

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

http://www.cas.org/infopolicy.html

=> s 14

L5

2 L4

=> d 1-2 bib fhitstr

L5 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2006 ACS on STN

AN 2005:116079 CAPLUS

DN 142:198489

TI Bridged carbocyclic compounds, their preparation, and use in polymerization

IN Van Court, Carr Richard; Markley, Thomas John; Abdourazak, Atteye Houssein; Marsella, John Anthony

PA Air Products and Chemicals, Inc., USA

SO Eur. Pat. Appl., 61 pp.

CODEN: EPXXDW

DT Patent

LA English

FAN.CNT 1

1111.011 1																			
	PATENT NO.			KIND DATE		}	APPLICATION NO.					DATE							
											- -								
ΡI	EP 1505050			A1 20050209		EP 2004-18304						20040802							
		R:	ΑT,	ΒE,	CH,	DE,	DK,	ES,	FR,	GB,	GR,	IT,	LI,	LU,	NL,	SE,	MC,	PT,	
			ΙE,	SI,	LT,	LV,	FI,	RO,	MK,	CY,	AL,	TR,	BG,	CZ,	EE,	HU,	PL,	SK,	HR
	US 2005037289 SG 108958			A1		2005	US 2004-784377				20040223 20040728								
				A1		20050228 SG 2004-4281													
	JP 2005053914			A2		2005	0303	JP 2004-228514				20040804							
PRAI	US	2003	-492	573P		P		2003	0804										
	US	2004	-784	377		Α		2004	0223										

OS MARPAT 142:198489

IT 838848-01-4P

RL: IMF (Industrial manufacture); PREP (Preparation)

(bridged carbocyclic compds., their preparation, and use in polymerization)

RN 838848-01-4 CAPLUS

CN 2-Hexanol, 5-(bicyclo[2.2.1]hept-5-en-2-yloxy)-1,1,1,6,6,6-hexafluoro-2,5-bis(trifluoromethyl)-, homopolymer (9CI) (CA INDEX NAME)

CM 1

CRN 838847-28-2 CMF C15 H14 F12 O2

RE.CNT 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

L5 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2006 ACS on STN

AN 2004:491231 CAPLUS

DN 142:490275

TI Novel reactions of quadricyclane. A new route to monomers for low-absorbing polymers in 157 nm photoresists

AU Marsella, John A.; Abdourazak, Atteye H.; Carr, Richard V. C.; Markley, Thomas J.; Robertson III, Eric A.

CS Corporate Science and Technology Center, Air Products and Chemicals, Inc., Allentown, PA, 18195, USA

SO Proceedings of SPIE-The International Society for Optical Engineering (2004), 5376(Pt. 1, Advances in Resist Technology and Processing XXI), 266-275

CODEN: PSISDG; ISSN: 0277-786X

PB SPIE-The International Society for Optical Engineering

DT Journal

LA English

IT 838848-01-4P

RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation) (preparation and properties of polymerized norbornene ethers obtained by reactions of quadricyclane with fluorinated diols for use in 157 nm photoresist formulations)

RN 838848-01-4 CAPLUS

CN 2-Hexanol, 5-(bicyclo[2.2.1]hept-5-en-2-yloxy)-1,1,1,6,6,6-hexafluoro-2,5-bis(trifluoromethyl)-, homopolymer (9CI) (CA INDEX NAME)

CM 1

CRN 838847-28-2 CMF C15 H14 F12 O2

RE.CNT 22 THERE ARE 22 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> file stnguide
COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 8.86 176.45

FULL ESTIMATED COST

FILE 'STNGUIDE' ENTERED AT 13:12:28 ON 15 APR 2006
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AND TECHNOLOGY CORPORATION, AND FACHINFORMATIONSZENTRUM KARLSRUHE

FILE CONTAINS CURRENT INFORMATION.

LAST RELOADED: Apr 7, 2006 (20060407/UP).

=> file req

COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 1.14 177.59

FULL ESTIMATED COST

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STRUCTURE FILE UPDATES: 13 APR 2006 HIGHEST RN 880388-58-9 DICTIONARY FILE UPDATES: 13 APR 2006 HIGHEST RN 880388-58-9

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TSCA INFORMATION NOW CURRENT THROUGH January 6, 2006

Please note that search-term pricing does apply when conducting SmartSELECT searches.

* The CA roles and document type information have been removed from * the IDE default display format and the ED field has been added, * effective March 20, 2005. A new display format, IDERL, is now * available and contains the CA role and document type information. *

Structure search iteration limits have been increased. See HELP SLIMITS for details.

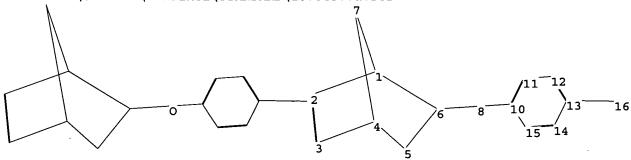
REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/ONLINE/UG/regprops.html

=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

Uploading C:\Documents and Settings\rkeys\My
Documents\STNEXP4\TEMPLATE\STANDARD\10784377a.str



```
chain nodes :
8 16
ring nodes :
1 2 3 4 5
             6 7 10
                       11
                           12
chain bonds :
6-8 8-10 13-16
ring bonds :
1-2 1-6 1-7 2-3
                       4-5
                            4-7
                                5-6 10-11 10-15 11-12 12-13 13-14 14-15
exact/norm bonds :
6-8 8-10
exact bonds :
1-2 1-6 1-7 2-3 3-4 4-5
normalized bonds :
10-11 10-15 11-12 12-13 13-14
isolated ring systems :
containing 1 : 10 :
```

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:CLASS 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:CLASS

L6 STRUCTURE UPLOADED

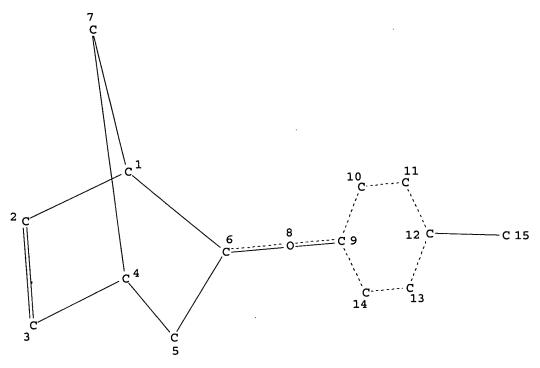
=> que L6

L7 QUE L6

=> d

L7 HAS NO ANSWERS

L6 ST



NODE ATTRIBUTES: NSPEC IS R AΤ NSPEC IS R AΤ NSPEC IS R ΑT NSPEC IS R ΑT NSPEC IS R ΑT NSPEC IS R ΑT 6 NSPEC IS R ΑT 7 NSPEC IS C ΑT 8 NSPEC IS R ΑT 9 NSPEC IS R AΤ 10 NSPEC IS R AΤ 11 NSPEC IS R AT 12 NSPEC IS R AΤ 13 NSPEC IS R AΤ 14 NSPEC IS C AΤ 15 DEFAULT MLEVEL IS ATOM MLEVEL IS CLASS AT DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RSPEC I

NUMBER OF NODES IS 15

STEREO ATTRIBUTES: NONE L7 QUE L6

=> s 17

SAMPLE SEARCH INITIATED 13:24:51 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED -90 TO ITERATE

100.0% PROCESSED

90 ITERATIONS

1 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE** BATCH **COMPLETE** PROJECTED ITERATIONS: 2369 1231 TO

PROJECTED ANSWERS:

1 TO 80

L8

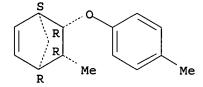
MF

1 SEA SSS SAM L6

=> d scan

REGISTRY COPYRIGHT 2006 ACS on STN L8 1 ANSWERS TN Bicyclo[2.2.1]hept-2-ene, 5-methyl-6-(4-methylphenoxy)-, (1R,4S,5S,6S)-rel- (9CI) C15 H18 O

Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> s 17 ful

FULL SEARCH INITIATED 13:25:07 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED -2147 TO ITERATE

100.0% PROCESSED 2147 ITERATIONS 12 ANSWERS

SEARCH TIME: 00.00.01

L9 12 SEA SSS FUL L6

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 167.38 344.97

FULL ESTIMATED COST

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L11

0 PHOTORESIST AND L10

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=> d his

(FILE 'HOME' ENTERED AT 13:08:37 ON 15 APR 2006)

FILE 'REGISTRY' ENTERED AT 13:08:53 ON 15 APR 2006

L1 STRUCTURE UPLOADED

L2 QUE L1

L3 1 S L2 L4 34 S L2 FUL

FILE 'CAPLUS' ENTERED AT 13:10:19 ON 15 APR 2006

L5 2 S L4

FILE 'STNGUIDE' ENTERED AT 13:12:28 ON 15 APR 2006

FILE 'REGISTRY' ENTERED AT 13:23:48 ON 15 APR 2006

L6 STRUCTURE UPLOADED

L7 QUE L6

L8 1 S L7

L9 12 S L7 FUL

FILE 'CAPLUS' ENTERED AT 13:25:14 ON 15 APR 2006

L10 7 S L9

L11 0 S PHOTORESIST AND L10

=> d l10 1-7 ti

- L10 ANSWER 1 OF 7 CAPLUS COPYRIGHT 2006 ACS on STN
- TI A two step, non-stereospecific cation radical Diels-Alder reaction
- L10 ANSWER 2 OF 7 CAPLUS COPYRIGHT 2006 ACS on STN
- TI Stereospecificity and Mechanism in Cation Radical Diels-Alder and Cyclobutanation Reactions
- L10 ANSWER 3 OF 7 CAPLUS COPYRIGHT 2006 ACS on STN
- TI 7,7-Dimethoxytetrachloronorbornenol ethers
- L10 ANSWER 4 OF 7 CAPLUS COPYRIGHT 2006 ACS on STN
- TI Vinyl ethers in diene synthesis. III. Synthesis of adducts of cyclopentadiene and hexachlorocyclopentadiene with vinyl ethers of phenols
- L10 ANSWER 5 OF 7 CAPLUS COPYRIGHT 2006 ACS on STN
- TI Simple vinyl ethers in diene synthesis. I. Reaction of vinyl ethers of phenol, cresols, and halophenols with hexachlorocyclopentadiene
- L10 ANSWER 6 OF 7 CAPLUS COPYRIGHT 2006 ACS on STN
- TI Vinyl compounds in Diels-Alder synthesis. III. Synthesis and properties of ethers of bicycloheptene series and di(endomethylene)octahydronaphthalene containing aromatic radicals
- L10 ANSWER 7 OF 7 CAPLUS COPYRIGHT 2006 ACS on STN
- TI Vinyl compounds in Diels-Alder reaction. Stereospecific orientation of Diels-Alder reaction of vinyl aryl ethers with cyclopentadiene as related to temperature

=> file caplus

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST ENTRY SESSION 6.10 351.07

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FILE COVERS 1907 - 15 Apr 2006 VOL 144 ISS 17 FILE LAST UPDATED: 14 Apr 2006 (20060414/ED)

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120 NORBON?

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36938 TRIFLUOROMETHYL

8663 CARBINOL

31 TRIFLUOROMETHYL CARBINOL

(TRIFLUOROMETHYL (W) CARBINOL)

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36938 TRIFLUOROMETHYL

8663 CARBINOL

31 TRIFLUOROMETHYL CARBINOL

(TRIFLUOROMETHYL (W) CARBINOL)

L13 3 NORBORN? AND (HEXFLUOROPROPANOL OR TRIFLUOROMETHYL CARBINOL)

=> d 1-3 ti

- L13 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN
- TI Equilibrium water uptake and diffusion behavior in model polynorbornene photoresist polymers
- L13 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN
- TI Effect of nanoscale confinement on the diffusion behavior of photoresist polymer thin films
- L13 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN
- TI Dissolution behavior of bis-trifluoromethyl-carbinol -substituted polynorbornenes

=> d 1-3 bib ab

- L13 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN
- AN 2005:592339 CAPLUS
- TI Equilibrium water uptake and diffusion behavior in model polynorbornene photoresist polymers
- AU Hoskins, Trevor; Roman, Paul J.; Ludovice, Peter J.; Henderson, Clifford L.

- CS Georgia Institute of Technology, School of Chemical & Biomolecular Engineering, Atlanta, GA, 30332-0100, USA
- SO Proceedings of SPIE-The International Society for Optical Engineering (2005), 5753(Pt. 2, Advances in Resist Technology and Processing XXII), 851-861

CODEN: PSISDG; ISSN: 0277-786X

- PB SPIE-The International Society for Optical Engineering
- DT Journal
- LA English
- AB As 193 nm immersion lithog. continues to evolve, the need to understand the effect of the immersing liquid on the resulting photoresist properties continues to grow. With this in mind, the sorption of water (using both liquid and vapor environments) in two model photoresist polymer resins based on functionalized poly norbornene) was examined using quartz crystal microbalance techniques. Similar to the results presented by Berger and coworkers, it was found that the water uptake in bistrifluoromethyl carbinol substituted polynorbornene (HFAPNB) increases as the polymer mol. weight increases, while the diffusion coefficient of water in these materials remains relatively constant over the

same

range in mol. weight In contrast, trifluorosulfonamide-substituted polynorbornene displays a relatively constant level of water uptake as a function of polymer mol. weight, while the diffusion coefficient decreases by

more

than an order of magnitude over the same mol. weight range. Sorption expts. performed as a function of temperature have shown that the water diffusion in these polynorbornene polymers can be described using an Arrhenius relationship. The activation energy of water diffusion was compared in both HFAPNB and poly(hydroxystyrene). The activation energy for diffusion of water in HFAPNB is substantially larger than in the case of poly(hydroxystyrene). This is consistent with the view that polynorbornenes possess relatively stiff and rigid backbones as compared to more flexible polymers such as poly(hydroxystyrene). The activation energy for water diffusion in HFAPNB was found to be a strong function of polymer mol. weight, with the activation energy decreasing with increasing mol. weight

RE.CNT 23 THERE ARE 23 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

L13 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN

AN 2004:491242 CAPLUS

DN 142:454199

- TI Effect of nanoscale confinement on the diffusion behavior of photoresist polymer thin films
- AU Singh, Lovejeet; Ludovice, Peter J.; Henderson, Clifford L.
- CS School of Chemical & Biomolecular Engineering, Georgia Institute of Technology, Atlanta, GA, 30332-0100, USA
- SO Proceedings of SPIE-The International Society for Optical Engineering (2004), 5376(Pt. 1, Advances in Resist Technology and Processing XXI), 369-378

CODEN: PSISDG; ISSN: 0277-786X

- PB SPIE-The International Society for Optical Engineering
- DT Journal
- LA English
- AB The influence of film thickness and mol. weight on the diffusion coeffs. of water, benzene, and trifluoroacetic acid in two photoresist polymers, poly(p-hydroxystyrene) and bis-trifluoromethyl carbinol substituted poly(norbornene), has been studied using quartz crystal microbalance (QCM) methods. Diffusion coeffs. for films as thin as approx. 50 nm were determined It was observed that the diffusion coefficient was a

strong function of film thickness, and that the diffusion coefficient decreases drastically as film thickness is reduced below a critical value. This critical

thickness value is found to be a function of both polymer structure and mol. weight In addition, the effect of film thickness on the equilibrium uptake of

the various penetrants was also determined In particular, the equilibrium water

uptake was shown to depend strongly on film thickness, polymer structure, and polymer mol. weight

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- L13 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN
- AN 2003:570067 CAPLUS
- DN 140:347337
- TI Dissolution behavior of bis-trifluoromethyl-carbinol -substituted polynorbornenes
- AU Hoskins, Trevor; Chung, Won Jae; Ludovice, Peter J.; Henderson, Clifford L.; Seger, Larry; Rhodes, Larry F.; Shick, Robert A.
- CS Georgia Institute of Technology, Atlanta, GA, 30332-0100, USA
- SO Proceedings of SPIE-The International Society for Optical Engineering (2003), 5039(Pt. 1, Advances in Resist Technology and Processing XX), 600-611
 - CODEN: PSISDG; ISSN: 0277-786X
- PB SPIE-The International Society for Optical Engineering
- DT Journal
- LA English
- AB As features shrink below 100 nm, new exposure technologies such as 157 nm lithog. are being developed. One of the critical challenges in developing these new lithog. tools and processes is the development of appropriate resist materials that can be used at these lower exposure wavelengths. Creating organic resist polymer resins for 157 nm exposure is a particularly challenging issue since many organic functional groups absorb at this wavelength. It has been previously shown that fluorinated polymers may offer the required low optical absorbance needed to serve as resist resins for 157 nm lithog. In particular, there has been interest in bistrifluoromethyl carbinol substituted polynorbornenes (HFAPNB) and similar materials for use in photoresists. The bistrifluoromethyl carbinol group offers a base soluble group that is sufficiently transparent to be used at 157 nm. This work has focused on the dissoln. behavior and other characteristics of bistrifluoromethyl carbinol substituted polynorbornenes. In particular, it was found that the dissoln. behavior of the HFAPNB homopolymer is strongly controlled by its ability to hydrogen bond with both neighboring chains and also other small mol. additives such as dissoln. inhibitors and photoacid generators. A detailed mol. level explanation for these effects is presented. The interaction of a series of com. photoacid generators with HFAPNB polymers are presented. The use of such information for the rational design of advanced resist materials using these polymers will be discussed.

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